Scheduling non-identical jobs on a batch resource

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Introduction

This paper discusses four different approaches, and several variations on them, to solving the problem of scheduling non-identical jobs on a batch processing machine. Batch processing machines, for the purposes of this paper, can process multiple, non-identical jobs simultaneously—but all jobs must be loaded into and unloaded from the machine at once, which introduces a considerable twist on the "simple parallel resources" known from typical example problems in existing literature.

The machines in question represents real-life resources like autoclaves or ovens, which can process multiple items at a time, but often cannot be opened at random—in fact, such machines often need to wait for the largest item in the batch to be done before the next batch can be inserted.

Malapert et al. [2012] proposed a global constraint programming (CP) algorithm consisting of a set of filtering rules to solve the problem. He achieved considerably better speeds than with a simple mixed-integer model (MIP), but it seems plausible that the new global constraint is unnecessarily cumbersome to achieve this performance; simple MIP, CP or decomposition approaches are easier to implement and extend. In this paper, we present 1) an improvement to his MIP model, 2) a CP model, 3) a decomposition approach to "divide and conquer" the problem, and 4) a new MIP model that greatly simplifies the original one.

1.1 Problem definition

We describe the problem as follows: assume we are given a set of jobs J, each of which has a processing time (or "length") p_j and a size (or "capacity requirement") s_j . Each job also has a due date d_j . The machine is characterized by its capacity b, and in every batch, the jobs' summed sizes must not exceed this number. All values are integer. 447723 The machine can only process one batch k of jobs at a time, and batches always take as long as the longest job in the batch (i.e. $P_k = \max_{j \in k}(p_j)$). Our objective is to minimize the lateness L of the latest job in J, where L is the difference between the job's completion time C_j and its due date d_j : in formal terms, min. $L_{max} = \max_j (C_j - d_j)$. The job's completion time, however, is the completion time of the batch, which in turn finishes with its *longest* job as stated above.

Malapert uses the standard format established by Graham et al. to summarize the problem as 1|p-batch;b| < n; non-identical $|L_{\max}|$, where p-batch;b| < n represents the parallel-batch nature and the finite capacity of the resource. A simpler version with identical job sizes was shown to be strongly NP-hard in [Brucker et al., 1998]; this problem, then, is no less difficult.

It helps to visualize the jobs before delving into the technicalities of scheduling them. Figure 1.1 shows a solution to a sample problem with eight jobs and a resource with capacity b = 20.

1.2 Organization of this paper

We first provide some background on the fundamental concepts of MIP and CP (sections ?? and ??), and review some of the most relevant publications on similar batch scheduling problems (section ??) before we describe Malapert's global constraint in some detail (section ??), as well as his original MIP model (section ??).

We then present possible improvements to the model in 4.2. Section 4.3 introduces a CP formulation of the same problem. Sections 4.4 and 4.4.1 describe a decomposition

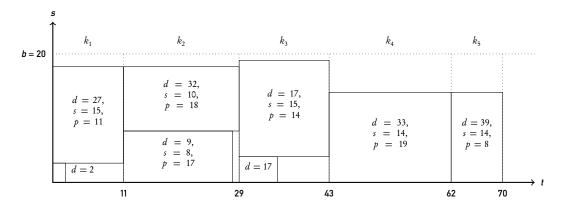


Figure 1.1: Optimal solution to an example problem with eight jobs (s_j and p_j not shown for the two small jobs)

approach. Finally, a new MIP approach is introduced in section ??.

An empirical comparison of the new models and a discussion of the results follow in sections 5.1 and 5.3. Ideas for future work are listed in 6.

Background

Many optimization problems, and scheduling problems in particular, are combinatorial in nature. The number of possible solutions grows exponentially with the number of input variables, and even with fast computers it is impossible to explore all of them individually to find the best one (also called "full enumeration", or "brute-force" search) in a reasonable amount of time. Often, however, it is possible to reason about subsets of solutions that are known to be suboptimal a priori. This limits the search space, allowing us to solve many instances of difficult combinatorial problems in few hours, minutes or even seconds.

Such constrained searches are often implemented in either one of two ways (or variants of them): as a *Constraint Programming model* (CP) or as a *Mixed Integer Programming model* (MIP). In this section I will briefly introduce the concepts behind both CP and MIP and review recent work on problems similar to the one dealt with here.

2.1 Constraint Programming

Many problems can be understood as a situation in which a set of values is to be chosen according to certain rules, but since the number of possible combinations of values is enormous, attempting to test all of them is infeasible.

Constraint programming (CP) is a formal framework for the formulation and solution of such problems. CP models consist of a set of variables $V = \{x_1, ..., x_n\}$, each x_i of which has a domain \mathcal{D}_i , the set of values that could conceivably be assigned to it. The CP problem is defined by a set of relationships that must hold between the variables. Model

2.1.1 illustrates the concept.

$$x \in \{1, \dots, 10\} \tag{2.1}$$

$$y \in \{9, \dots, 20\} \tag{2.2}$$

$$z \in \{1, \dots, 50\} \tag{2.3}$$

$$x > y \tag{2.4}$$

$$z \geqslant xy - 42 \tag{2.5}$$

Model 2.1.1: A simple CP model

Here, x, y, z are the variables with their respective domains.

2.1.1 Propagation

We can choose a variable and a value to assign to it, and explore the consequences of this assignment. Assign x = 5, then based on the constraint x > y only values < 5 are permissible for y, but none are available in \mathcal{D}_y . The assignment x = 5 leads to an *inconsistency*, and another value must be chosen.

If we choose x = 10 (which, clearly, is the only feasible value for x, since all others violate x > y), y is limited to y = 9. This means that $z \ge 9 \cdot 10 - 42 = 48$, thus $z = \{48, 49, 50\}$.

This process of successive elimination is known as *propagation*, and it is the core of all CP solver algorithms. In propagation, a variable's domain is reduced, and the constraints are used to reduce the domains of other variables accordingly, until no constraints are violated.

2.1.2 Consistency

Propagation is commonly seen as a way to enforce consistency. It can be triggered by fixing a variable to a value, but also by systematic evaluation of constraints. Let any two variables in a problem be involved in a binary constraint C, such as x > y above. If their domains are reduced such that they fulfill C (that is, $\forall a \exists b$ such that $a \in \mathcal{D}_x$, $b \in \mathcal{D}_y$, C(x = a, y = a,

b) = true), they are *arc-consistent* with regards to C. Propagation will commonly enforce arc consistency across all variables and constraints.

Arc consistency alone is not sufficient to guarantee the existence of a solution (satisfiability), unless the graph of variables and constraints is acyclic. Consider the following problem (Model 2.1.2):

$$x_1 = x_2 \tag{2.6}$$

$$x_1 = x_3 \tag{2.7}$$

$$x_2 = x_4 \tag{2.8}$$

$$x_3 \neq x_4 \tag{2.9}$$

$$x_1, x_2, x_3, x_4 \in \{1, 2\}$$
 (2.10)

Model 2.1.2: Unsatisfiable CP problem

The model is arc-consistent but has no solution: forcing any variable to assume a fixed value will result in at least constraint being violated.

The notion of arc consistency (two variables) can be extended, e.g. to three variables related through binary constraints (*path consistency*) or to constraints involving more than two variables, where every value assigned to a variable *x* must be consistent with all of all other variables' values (*generalized arc consistency*).

2.1.3 Global constraints

Propagation makes use of graph-theory based algorithms that enforce a form of consistency on the variables involved in a constraint. Most constraints are formulated in the form of binary relationships involving a single operator (e.g. \geqslant ,= or \neq). But it is often possible and beneficial to impose constraints on groups of variables at once. Consider the following classic example: This model prescribes that the variables x_1, x_2, x_3, x_4 are to take on different values. The model is arc consistent. To find a solution (or insatisfiability), a value has to be assigned to three of the variables, and the solver has to propagate the domain reductions after every assignment. Consider, on the other hand, the use of a specialized

$$x_{1} \neq x_{2}$$
 (2.11)

$$x_{2} \neq x_{3}$$
 (2.12)

$$x_{3} \neq x_{4}$$
 (2.13)

$$x_{4} \neq x_{1}$$
 (2.14)

$$x_{1} \neq x_{3}$$
 (2.15)

$$x_{2} \neq x_{4}$$
 (2.16)

$$x_{1}, x_{2}, x_{3}, x_{4} \in \{1, 2, 3\}$$
 (2.17)

Model 2.1.3: A clique of not-equal constraints

constraint called alldifferent(x_1, x_2, x_3, x_4). This constraint, implemented as a custom algorithm, can speed up the propagation rapidly. In the case of our example, the constraint can infer that the number of variables exceeds the number of available values. The solver can conclude that the problem is insatisfiable without any propagation.

Such specialized or "global" constraints can efficiently exploit properties that hold for certain relationships among groups of variables, and that would not be available if the relationship were expressed as a set of binary constraints.

2.2 Mixed Integer Programming

2.2.1 Linear Programming

Mixed Integer Programs (or "MIP models") express the minimization of a linear function subject to linear constraints. If all variables in the problem can be rational in the solution, the MIP model is really a *linear program* (LP), which can be solved in polynomial time.¹

Figure 2.1 illustrates the concept of an LP in two variables, x_1 and x_2 , as listed in model 2.2.1. The set of feasible solutions is given by the shaded area bounded by the axes and by three inequalities ("constraints"). A third linear term, the objective function $3x_1 + x_2$, is to

¹In practice, variations on Dantzig's *simplex method* are most often used to solve LPs. Such solvers perform very well on most problems, but no known variant has been proven to have polynomial worst-case complexity [?]. Solvers with theoretically polynomial-time complexity exist (Karmarkar's algorithm Karmarkar [1984] has a runtime of $\mathcal{O}(n^{3.5}L^2 \cdot \log L \cdot \log \log L)$, for instance, where L is the number of bits of input), but are used less frequently.

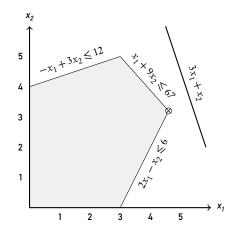


Figure 2.1: Graphical representation of a typical LP problem

be maximized.²

Maximize
$$3x_1 + x_2$$
 (2.18)
subject to the constraints $-x_1 + 3x_2 \le 12$ (2.19)
 $x_1 + 9x_2 \le 67$ (2.20)
 $2x_1 - x_2 \le 6$ (2.21)
 $x_1 \ge 0$ (2.22)
 $x_2 \ge 0$ (2.23)

Model 2.2.1: A simple LP model, as shown in figure 2.1

Although the objective function is shown in the figure as a line in a specific location, note that, as it is not an equation, it can be represented by any line parallel to that shown in the figure. While we are trying to maximize the objective value, the solution must be feasible. It is thus obvious that the desired extreme value of our objective function is found at one of the "corners" of the shaded area: the solution is marked \otimes in the figure. In fact, since the shaded area generated by linear inequalities will always be a convex polygon (or polyhedron, in higher dimensions), the solution will invariably be found at an intersection of hyperplanes.

²Minimization is more common, but note that multiplying the objective by -1 achieves this.

2.2.2 Solving MIP models using branch-and-bound

MIP models are LPs in which some of the decision variables are declared integers—thus their name. Like LP models, MIP models also require an objective function to be minimized. Figure 2.2a illustrates this based on the LP problem above: now, only the black dots

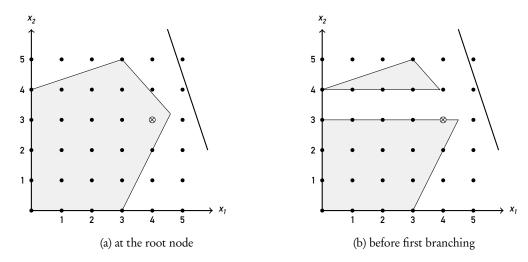


Figure 2.2: Graphical representation of a typical MIP problem

represent feasible solutions. While the solution is easily found in the figure by inspection (simply round to the nearest feasible integral solution!), this is not the case in problems with many variables; it is difficult enough to visualize the problem in three dimensions, and many problems require hundreds or thousands. Moreover, rounding is particularly unreliable with variables of small domains (e.g. binary variables), which are often used in MIP models to represent decisions. Indeed, solving MIP models is NP-hard.

Similar to CP searches, MIP searches can be thought of as trees, where every branch represents an assignment of a value to a domain and every leaf represents a feasible solution. The difference lies in the way MIP explores this search tree.

The MIP solver first solves the problem as an LP, which will usually result in fractional values for all or most of the variables. In this context, the LP is known as the *LP relaxation* of the problem—an easier, approximate version of the original. Assuming we are minimizing the objective, the resulting LP objective value will be a lower bound on the optimal

MIP objective value. At this point, the solution is $x_1 = 4.6$, $x_2 = 3.2$.

The solver then chooses a variable, based on heuristics, and *branches* on it. In this case, assume that we branch on x_2 , which means that we set $x_2 \le 3.0 \lor x_2 \le 4.0$. We now have two new MIPs, as shown in figure 2.2b. Both of these new subproblems are need solving.³ At this point, another heuristic decides which subproblem is solved first. Again, the chosen subproblem is first solved in its LP relaxation, and the above procedure is repeated until an integral solution is found. If this solution is the best solution known so far, it is stored as the *incumbent*. The solver then *backtracks*: it undoes some of the previously fixed variables, and explores other (previously unsolved) subproblems.

This behaviour can be visualized as a search tree in which every node represents a decision between two subproblems. At every node, a subproblem is chosen, the subproblem's LP relaxation is run, and the next variable is chosen to be branched on.

When the solver backtracks to a previous node to explore another subproblem, say P_1 , it can compare the LP solution of P_1 with the current incumbent's solution. Since a solution will never improve with additional integrality constraints, the subtree of P_1 can be ignored if the objective value of the LP solution to P_1 is worse than that of the current incumbent. This "pruning" of the search tree during the search is referred to as *bounding*, and is what allows many MIP models to be solved in less time than their theoretically exponential complexity suggests.

2.2.3 Branch-and-cut methods

2.2.4 Special ordered sets

2.2.5 Lazy constraints in MIP

Achieving fast search times in MIP requires striking a delicate balance between keeping the model small (reducing the time needed to solve the LP relaxations) and formulating the model tightly (pruning more branches off the search tree). Often, the way certain

³In this example, the optimal value for x_2 was within ± 1 of its LP value. This is often the case, and it is not immediately obvious from the figure why fixing $x_2 = 3.0 \lor x_2 = 4.0$ is insufficient. In some problems, however, the shaded polyhedron protrudes relatively far diagonally through the grid of integer solutions, in which case branching on $\leq \lfloor x_i \rfloor$ and $\geq \lceil x_i \rceil$ is necessary to capture the optimal feasible solution.

constraints can be used to perform pre-solve reductions plays into the performance also, and is difficult to predict.

Lazy constraints are one way to avoid this problem. They are expressed as regular linear constraints and checked against whenever the model encounters a new solution. Lazy constraints that were violated are then added to the model to be used in future instances.

This is particularly useful when a problem lends itself to the generation of vast sets of constraints involving few variables, but most of these constraints are likely not needed. In such cases, lazy constraints can greatly speed up solution times.

2.3 Scheduling using CP and MIP

- 2.3.1 Common types of scheduling problems
- 2.3.2 Temporal problems in CP
- 2.3.3 Temporal problems in MIP

2.4 Literature

The problem at hand is based on the work of Malapert et al. [2012], who proposed a global constraint sequenceEDD to be used in combination with pack to solve it to optimality. The sequenceEDD constraint is implemented as four distinct filtering rules applied at relevant domain changes: three to update bounds on $L_{\rm max}$ based on different conditions, and one to limit the number of batches based on the marginal cost difference between adding a job to an empty batch vs. an existing one. The new global was shown to significantly outperform a simple MIP model of the same problem.

Other authors have examined similar problems: Azizoglu and Webster [2000] provide an exact method and a heuristic for the same problem, but minimize makespan (C_{max}) instead of L_{max} , as have Dupont and Dhaenens-Flipo [2002]. Similar exact methods have been proposed for multi-agent variants with different objective functions [Sabouni and Jolai, 2010], for makespan minimization on single batch machines [Kashan et al., 2009], and

for makespan minimization on parallel batch machines with different release dates [Ozturk et al., 2012]. A more extensive review of MIP model applications in batch processing is given by Grossmann [1992].

State of the Art: Recent work by Malapert

3.1 The sequenceEDD global constraint

3.2 Malapert's MIP formulation

Malapert's original MIP approach is given in Model 3.2.1. It uses a set of binary decision variables x_{jk} to represent whether job j is assigned to batch k. The original model assumes a set K of |K| = |J| batches, the number of jobs being a trivial upper bound on the number of batches required; it also enforces an earliest-due-date-first (EDD) ordering of the batches (constraint (3.7)).

Constraints (3.2) ensure that every job j is assigned to one batch k, and one batch only. Constraints (3.3) ensure that no batch exceeds the machine's capacity b. Constraints (3.4) define each batch's processing time P_k as the maximum processing time of the jobs j assinged to it. Constraints (3.5) define each batch's completion time C_k as that of the previous batch, plus the previous batch's processing time. Constraints (3.7) ensure that batches are ordered by due date. Constraints (3.8) define the objective value L_{\max} .

Model 3.2.1: Malapert's original MIP model

Modelling the problem

In this section we present the models examined. Beginning with the MIP model used by Malapert, we explore several additional constraints that, while redundant, tighten the search space. We comment on the performance of the approaches (see section ?? for a comprehensive comparison of performance). We present a CP model and a decomposition-based approach using either MIP or CP. Finally, we introduce a new approach to setting up a MIP model that significantly improves upon the original MIP and outperforms all other models presented in this paper.

All MIP and CP models were implemented in C++ using IBM Ilog's Cplex and CP Optimizer solvers, respectively.

4.1 MIP model

We first replicated Malapert's original MIP approach, as given in Model 3.2.1, to establish a performance baseline to compare other models to. The original model assumes a set K of |K| = |J| batches, the number of jobs being a trivial upper bound on the number of batches required; it also enforces an earliest-due-date-first (EDD) ordering of the batches (constraint (3.7) in Model 3.2.1).

4.2 Improved MIP model

Several improvements can be made to Malapert's MIP model in the form of additional constraints shown in Model 4.2.1, as described in greater detail in the subsections below.

$$\sum_{j \in J} x_{j,k-1} = 0 \longrightarrow \sum_{j \in J} x_{jk} = 0 \quad \forall k \in K$$

$$(4.1)$$

$$e_k + \sum_{j \in J} x_{jk} \ge 1 \qquad \forall k \in K$$

$$n_j(e_k - 1) + \sum_{j \in J} x_{jk} \le 0 \qquad \forall k \in K$$

$$(4.2)$$

$$n_j(e_k - 1) + \sum_{i \in I} x_{jk} \le 0 \qquad \forall k \in K$$

$$(4.3)$$

$$e_k - e_{k-1} \geqslant 0 \qquad \forall k \in K \tag{4.4}$$

$$x_{ik} = 0 \qquad \qquad \forall \{j \in J, k \in K | j > k\} \tag{4.5}$$

$$e_{k} - e_{k-1} \geqslant 0 \qquad \forall k \in K \qquad (4.4)$$

$$x_{jk} = 0 \qquad \forall \{j \in J, k \in K | j > k\} \qquad (4.5)$$

$$L_{\max} \geqslant \lceil \frac{1}{b} \sum_{j} s_{j} p_{j} \rceil - \delta_{q} \qquad \forall q, \forall \{j \in J | d_{j} \leqslant \delta_{q}\} \qquad (4.6)$$

Model 4.2.1: Improvements to Malapert's original MIP model

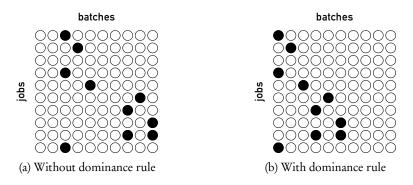


Figure 4.1: Dominance rule to eliminate empty batches followed by non-empty batches (circles represent the x_{jk} variables; a filled circle stands for $x_{jk} = 1$)

Grouping empty batches 4.2.1

The given formulation lacks a rule that ensures that no empty batch is followed by a nonempty batch. Empty batches have no processing time and a due date only bounded by d_{max} , so they can be sequenced between non-empty batches without negatively affecting $L_{\rm max}$. Since, however, desirable schedules have no empty batches scattered throughout, we can easily reduce the search space by disallowing such arrangements. The idea is illustrated in Figure 4.1.

A mathematical formulation is

$$\sum_{j \in J} x_{j,k-1} = 0 \longrightarrow \sum_{j \in J} x_{jk} = 0 \quad \forall k \in K.$$

$$(4.7)$$

To implement this, we can write constraints in terms of an additional set of binary variables, e_k , indicating whether a batch k is empty or not:

$$e_k + \sum_{j \in J} x_{jk} \geqslant 1 \qquad \forall k \in K,$$
 (4.8)

$$n_{j}(e_{k}-1)+\sum_{j\in I}x_{jk}\leqslant 0 \quad \forall k\in K. \tag{4.9}$$

Constraints (4.8) enforce $e_k = 1$ when the batch k is empty. Constraints (4.9) enforce $e_k = 0$ otherwise, since the sum term will never exceed the number of jobs n_j . The rule (4.7) can now be expressed as $e_{k-1} = 1 \rightarrow e_k = 1$, and implemented as follows:

$$e_k - e_{k-1} \geqslant 0 \quad \forall k \in K. \tag{4.10}$$

We can also prune any attempts to leave the first batch empty by adding a constraint $e_0 = 0$.

4.2.2 No postponing of jobs to later batches

Consider a schedule in which all jobs are assigned to a single batch (i.e. $x_{jk} = 1 \forall j = k$), ordered by non-decreasing due date, and by non-decreasing processing time in case of a tie between two jobs. For the purposes of this paper, we refer to this schedule as the *single-EDD* schedule.

Since, as shown above, there exists an optimal solution with EDD-ordered batches, we can restrict the search to such solutions. Any solution in which a job j is moved into a batch h later than its single-EDD batch k, i.e. h > j = k, necessarily results in a non-EDD solution. We can therefore write

$$x_{jk} = 0 \quad \forall \{j \in J, k \in K | j > k\}$$
 (4.11)

to exclude solutions in which jobs are assigned to later batches than their respective single-EDD batch.

4.2.3 Lower bound on L_{max}

Let a bucket q denote the set of all batches with due date δ_q . Then the completion date C_q of this bucket is the completion date of the last-scheduled batch with due date δ_q , and the lateness of the bucket q is $L_q = C_q - \delta_q$. Since all batches up to and including those in bucket q are guaranteed to contain all jobs with due dates $d \leq \delta_q$, as ensured by the EDD ordering of batches, the lower bound on every bucket's lateness $LB(L_q)$ is a valid lower bound on L_{\max} . In other words, jobs with due date $d \leq \delta_q$ will be found only in batches up to and including the last batch of bucket q. This provides a lower bound on the lateness of bucket q:

$$L_{\max} \geqslant C_{\max,q} - \delta_q \quad \forall q \tag{4.12}$$

The buckets up to bucket q will likely also contain some later $(d > \delta_q)$ jobs in the optimal solution, which weakens $LB(L_q)$'s usefulness but does not affect its validity as a lower bound.

Now we need to find $C_{\max,q}$, or at least a lower bound on it, in polynomial time. The simplest approach simply considers the jobs' total "area" (or "energy"), i.e. the sum of all $s_i p_i$ products:

$$C_{\max,q} \geqslant \lceil \frac{1}{b} \sum_{j} s_{j} p_{j} \rceil \quad \forall q, \forall \{j \in J | d_{j} \leqslant \delta_{q} \}$$
 (4.13)

A better lower bound on $C_{\max,q}$ would be given by a preemptive-cumulative schedule. Unfortunately, minimizing C_{\max} for such problems is equivalent to solving a standard bin-packing problem, which requires exponential time.¹

¹In a preemptive-cumulative schedule, jobs may be stopped and restarted mid-execution, but occupy a constant amount s_j on the resource while executing. In such a schedule, minimizing the makespan is as difficult as solving a bin-packing problem: we can break jobs into small pieces (no longer than the smallest common divisor of the jobs' lengths p) and then pack them together such as to minimize the number of small time slots needed.

4.3 CP model

The constraint programming model is based on a set of decision variables $B_j = \{k_1, ..., k_{n_k}\}$, where each variable stands for the batch job j is assigned to. The complete model is given in Model 4.3.1.

$$Min. L_{max} (4.14)$$

s.t.
$$\operatorname{pack}(J, K, b)$$
 (4.15)

$$cumul(J,b) (4.16)$$

$$P_k = \max_j p_j \qquad \qquad \forall \{j \in J | B_j = k\}, \forall k \in K$$
 (4.17)

$$D_k = \min_j d_j \qquad \qquad \forall \{j \in J | B_j = k\}, \forall k \in K$$
 (4.18)

$$C_k + P_{k+1} = C_{k+1} \qquad \forall k \in K \tag{4.19}$$

$$L_{\max} \geqslant \max_{k} \left(C_k - D_k \right) \tag{4.20}$$

$$B_{j} \leq k \qquad \qquad \forall \{j \in J, k \in K | j > k\} \tag{4.22}$$

Model 4.3.1: Constraint programming model

Constraint (4.15) makes sure the jobs are distributed into the batches such that no batch exceeds the capacity b. Constraint (4.16), a global cumulative constraint, keeps jobs from overlapping on the resource—while redundant with (4.15), this speeds up propagation slightly. Constraints (4.17), (4.18), (4.19) and (4.20) define P_k , D_k , C_k (batch completion time) and L_{\max} , respectively.

4.3.1 Grouping empty batches

We can force empty batches to the back and thus establish dominance of certain solutions. The implementation is much easier than in the MIP model:

If Then
$$(P_k = 0, P_{k+1} = 0) \quad \forall k \in \{k_1, \dots, k_{n_k-1}\}$$
 (4.23)

4.3.2 No postponing of jobs to later batches

Just like in the MIP model, jobs should never go into a batch with an index greater than their own:

$$B_{j} \leq k \quad \forall \{j \in J, k \in K | j > k\} \tag{4.24}$$

4.4 Decomposition approach

Instead of solving the entire problem using one model, we can solve the problem step by step, using the best techniques available for each subproblem. This approach is inspired by *Benders decomposition*.

A basic version of this approach uses branch-and-bound to transverse the search tree. At each node on level ℓ , a single MIP and/or CP model is run to assign jobs to batch $k=\ell$. The remaining jobs are passed to the children nodes, which assign jobs to the next batch, and so on—until a solution, and thus a new upper bound on $L_{\rm max}$, is found. Several constraints are used to prune parts of the search tree that are known to offer only solutions worse than this upper bound. Figure 4.2 shows an example in which a MIP model is used to assign jobs to the batch at every node. Algorithm 1 outlines what happens at each node:

Algorithm 1 MIP node class code overview

```
update currentAssignments
                                                                                     ▶ this keeps track of where we are in the tree
if no jobs given to this node then
                                                                        ▶ if this is a leaf node, i.e. all jobs are assigned to batches
    calculate L_{\text{max,current}} based on \textit{currentAssignments}
    if L_{\text{max,current}} < L_{\text{max,incumbent}} then
        L_{\text{max,incumbent}} \leftarrow L_{\text{max,current}}
        bestAssignments \leftarrow currentAssignments
    return to parent node
end if
set up MIP model
                                                                                                                 ▶ as described below
repeat
                                                                                                ▶ let model assign jobs to the batch
    x_i \leftarrow \text{model.solve}(x_i)
    spawn and run child node with all \{j | x_i = 0\}
                                                                                                  ▶ pass unassigned jobs to children
    add constraint to keep this solution from recurring
                                                                                       ▶ this happens once the child node returns
until model has no more solutions
return to parent node
```

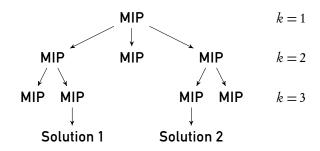


Figure 4.2: Batch-by-batch decomposition using MIP only

the model finds the best jobs to assign to the batch according to some rule, lets the children handle the remaining jobs, and tries the next best solution once the first child has explored its subtree and backtracked.

Using MIP and cumulative packing after the batch

The first version of the branch-and-bound batch-by-batch decomposition uses a MIP model at each node to assign jobs to the respective batch. The remaining jobs are packed such as to minimize their $L_{\rm max}$, with a relaxation of the batching requirement, i.e., as if on a cumulative resource.

x_{j}	is 1 iff job j is assigned to the batch
u_{jt}	is 1 iff job j starts in time slot t
T_{jt}	is the set of all time slots occupied by job j if it ended at time t , that is
, -	$T_{it} = \{t - p_i + 1, \dots, t\}$
v_k	is the start time of the batch at the given node in the search tree
$L_{\rm max,incmb}$	is the incumbent (known best) value of and thus an upper bound on $L_{\rm max}$
\mathcal{H}	is the set of all indexed time points $\{0, \dots, n_t - 1\}$
w_{i}	is 1 iff job j is either longer than the batch $(p_j > P_k)$ or already part of the
,	batch ($x_i = 1$). Neither condition must be fulfilled for constraint (4.36) to
	have an effect on job j

Table 4.1: Notation used in the decomposition model

Model 4.4.1 implements a time-indexed cumulative constraint on the non-batched jobs. Constraints (4.26) through (4.28) ensure that the batch stays below capacity, define the

Min.
$$L_{\text{max,cumul}}$$
 (4.25)

s. t.
$$\sum_{j} s_{j} x_{j} \leq b$$
 $\forall j \in J$ (4.26)

$$P_k \geqslant p_j x_j \qquad \forall j \in J \tag{4.27}$$

$$\sum_{j} x_{j} \geqslant 1 \qquad \qquad \forall \{j \in J | d_{j} = \min(d_{j})\}$$
 (4.28)

$$P_k + \frac{1}{b} \sum_i s_i p_i \leqslant d_j + L_{\text{max,incmb}} - 1 - v_k \quad \forall j \in J, \forall \{i \in J | d_i \leqslant d_j\}$$
 (4.29)

$$\sum u_{jt} = 1 \qquad \forall j \in J \tag{4.30}$$

$$\sum_{j} \sum_{t' \in T_{i,t}} s_j u_{jt'} \leqslant b \qquad \forall t \in \mathcal{H}$$
 (4.31)

$$(v_k + t + p_j)u_{jt} \le d_j + L_{\text{max.incmb}} - 1 \qquad \forall j \in J, \forall t \in \mathcal{H}$$

$$(4.32)$$

$$L_{\text{max,cumul}} \ge (v_k + t + p_j)u_{jt} - d_j \qquad \forall j \in J, \forall t \in \mathcal{H}$$
(4.33)

$$u_{j,t=0} = x_j \tag{4.34}$$

$$u_{it} \le (1 - x_j)$$
 $\forall i, j \in J, \forall t \in \{1, ..., p_j - 1\}$ (4.35)

$$b - \sum_{i \in I} s_i x_i \le (b \, w_j + 1) s_j \qquad \forall j \in J$$
 (4.36)

$$P_k + 2w_j n_t \geqslant p_j + n_t x_j \qquad \forall j \in J \tag{4.37}$$

$$P_k - 2(1 - w_j)n_t \leqslant p_j + n_t x_j - 1 \qquad \forall j \in J$$
 (4.38)

Model 4.4.1: MIP model in batch-by-batch branch-and-bound

duration of the batch P_k and force at least one of the earliest-due jobs into the batch.

Constraints (4.29) express the interval relaxation used to ensure that no jobs exceed their latest allowable finish date given any batch assignment. Even jobs that are assigned to the batch have to fulfill this requirement.

Constraints (4.30) and (4.31) implement the cumulative nature of the post-batch assignments by ensuring that each job starts only once, and no jobs overlap on a given resource at any time.

Constraints (4.32) again limit the possible end dates of a job, but unlike (4.29), they use the time assignments on the cumulative resource to determine end dates. Constraints (4.33) define the value of $L_{\text{max,cumul}}$, the maximum lateness of any job in the non-batched set.

Constraints (4.34) force batched jobs to start at t = 0, while (4.35) force *all* jobs to start either at t = 0 or after the last batched job ends.

Constraints (4.36) enforce a dominance rule: jobs must be assigned to the batch such that the remaining capacity, $b_r = b - \sum_j s_j x_j$, is less than the size s_j of the *smallest* job from the set of non-batched jobs that are *not longer* than the current batch. That is, if there exists an non-batched job j with $p_j \leq P_k$ and $s_j \leq b_r$, then the current assignment of jobs is infeasible in the model. The reasoning goes as follows: given any feasible schedule, assume there is a batch k_a with b_r remaining capacity and a later batch k_b containing a job j such that $s_j \leq b_r$ and $p_j \leq P_{k_a}$. Then job j can always be moved to batch k_a without negatively affecting the quality of the solution: if the schedule was optimal, then moving j will not affect L_{\max} at all (otherwise, it was no optimal schedule); if the schedule was not optimal, then L_{\max} will stay constant (unless j was the longest job in k_b and L_{\max} occured in or in a batch after k_b , in which case L_{\max} will be improved).

This rule is implemented by means of a binary variable w_j , which, as defined by constraints (4.36) and (4.37), is 1 iff $p_j > P_k \lor x_j = 1$. These are the cases in which a job j is not to be considered in (4.35), and so w_j is used to scale s_j to a value insignificantly large in the eyes of the constraint's less-than relation.

After a solution is found, a child node in the search tree has run the subtree and returned, a constraint of the form

$$\sum_{j} x_{j} + \sum_{i} (1 - x_{i}) \le n_{j} - 1 \quad \forall \{j \in J | x_{j} = 1\}, \forall \{i \in J | x_{i} = 0\}$$
 (4.39)

is added to the model before the solver is called again, to exclude the last solution from the set of feasible solutions.

4.4.1 Using CP and cumulative packing after the batch

This approach is equivalent, but we now use CP to select the batch assignments, again based on a minimized L_{max} among the non-batched jobs. Model 4.4.2 uses interval variables j to represent the jobs; time constraints on the jobs ("est", "lft" etc.) are represented as functions such as start0f(j) and end0f(j).

Contraints (4.41) through (4.44) bi-directionally define the relationship between a job's x_j and start0f(j): $x_j = 1$ is equivalent with a start time of t = 0, and $x_j = 0$ is equivalent with a start time of $t \ge P_k$. The term $\sum_{i \in J} p_i$ is used as a large constant as it is greater than any job's start time.

Constraint (4.48) is equivalent to constraints (4.36) through (4.38) in Model 4.4.1 above. The cumulative constraint (4.51) ensures that jobs do not overlap.

4.5 Move-based MIP model

This approach is based on the idea of preserving EDD ordering among the batches. Let an initial schedule be the single-EDD schedule as defined in section 4.2.2. We can calculate the lateness $L_{k,\text{single}}$ of every job in this schedule in linear time.

For the purpose of this discussion, we use the following terminology: in any batch k holding multiple jobs in an EDD schedule, let *host job* denote the earliest-due job in the batch, and *guest jobs* all other jobs. If a batch k only holds one job j, then job j is said to be *single*.

We now observe the following two things:

Proposition 1. Consider a schedule in which batches are ordered by EDD. Then moving job j from batch k_{β} into an earlier batch k_{α} will preserve EDD ordering if j is single in k_{β} and if k_{α} is not empty.

Proof. Moving j into k_{α} will leave D_{α} unaffected, since k_{α} is not empty, i.e. there is a host job in k_{α} . Since the original schedule was EDD-ordered, the host in k_{α} is due earlier than j. Note that if k_{α} is empty, the move will potentially disrupt EDD ordering, and is thus

$$Min. L_{max} (4.40)$$

$$\text{s.t. } \operatorname{startOf}(j) \geqslant (1-x_j)P_k \qquad \qquad \forall j \in J \tag{4.41}$$

$$\mathsf{startOf}(j) \leqslant (1 - x_j) \sum_{i \in I} p_i \qquad \forall j \in J$$
 (4.42)

$$x_{j} \geqslant \frac{\frac{1}{2} - \operatorname{start0f}(j)}{\sum_{i \in J} p_{i}} \qquad \forall j \in J$$
 (4.43)

$$x_{j} \leqslant 1 + \frac{\frac{1}{2} - \operatorname{startOf}(j)}{\sum_{i \in J} p_{i}}$$
 $\forall j \in J$ (4.44)

$$P_k \geqslant \max_{i \in I}(x_i \, p_i) \tag{4.45}$$

$$endOf(j) = d_j + L_{max,incmb} - 1 \forall j \in J (4.46)$$

$$L_{\max} \geqslant \max_{j \in I} (\operatorname{endOf}(j) - d_j) \tag{4.47}$$

$$P_k + \frac{1}{b} \sum_i s_i p_i \leqslant d_j + L_{\text{max,incmb}} - 1 - v_k \qquad \forall j \in J, \forall \{i \in J | d_i \leqslant d_j\}$$
 (4.49)

$$\sum_{j} x_{j} \geqslant 1 \qquad \qquad \forall \{j \in J | d_{j} = \min(d_{j})\} \tag{4.50}$$

$$\operatorname{cumul}(J,b) \tag{4.51}$$

Model 4.4.2: CP model in batch-by-batch branch-and-bound

prohibited.

Batch k_{β} will be reduced to zero processing time, but batches after k_{β} will still be due after $k_{\beta-1}$, so EDD ordering is preserved.

Proposition 2. Starting from a single-EDD schedule, moving single jobs back into earlier non-empty batches will generate all possible EDD schedules, with the exception of those EDD schedules in which two hosts of the same due date are sorted in order of decreasing processing time.

Proof. The order in which moves are performed is arbitrary; moves are independent of

each other (capacity requirements notwithstanding). Therefore, any job can be moved into any earlier bucket (i.e. set of batches of equal due date). Within buckets, jobs are free to move with the only condition that host jobs be sorted by non-decreasing processing time, as given by the intial single-EDD schedule. Under this condition, then, all possible EDD schedules can be generated by moving single jobs into earlier batches.

Consider now a single-EDD schedule. Moving j from its single-EDD batch k_j into an earlier batch k_{α} has the following effect:

- the lateness of all batches after k_i is reduced by p_i ,
- the lateness of all batches after k_{α} , including those after k_{j} , is increased by max(0, $p_{j}-P_{\alpha}$), where P_{α} is the processing time of batch α before j is moved into it.

Since, in any batch k, only the host job's lateness (with index j=k) is relevant to L_{\max} , we can understand the lateness of batch k as the single-EDD lateness of job $k=L_{k,\text{single}}$, modified by summed effect all moves of other jobs into and out of earlier batches have on k, as listed above.

The following expression defines the lateness of a batch *k* based on this calculation:

$$L_{k} = L_{k,\text{single}} - \sum_{b=0}^{k} P_{b} - p_{b}(2 - x_{bb}) \quad \forall k \in K^{*}$$
 (4.52)

where x_{jk} is 1 iff job j is assigned to batch k in the solution schedule. The sum term represents the effects that all moves have on previous batches up to, and including, k: $x_{bh} = 1$ iff job h is not moved from its single-EDD batch h, and so the sum term evaluates to $P_h - p_h$ for a batch in which the host was not moved. If h has guests, then $P_h - p_h$ is positive if any of the guests is longer than j_h (i.e. $P_h > p_h$); if not, or if h has no guests at all, it will be zero. Effectively thus, for batch h, the time $\max(0, P_h - p_h)$ is added to L_k . If, however, job h was moved out of its batch, then the batch will have a processing time of $P_h = p_h$ (the fixed lower bound for P_h). In this case, effectively, for batch h, the time p_h is subtracted from L_k . The net sum of these additions and subtractions to and from $L_{k,\text{single}}$

adjusts the lateness of batch k to its correct number given the values of x_{ik} .

Note also that we only need to calculate the lateness for $k \in K^*$, the set of batches that are the last in their respective buckets; in other words, those with the longest processing time in their respective bucket. This is true because empty batches in a bucket will be reduced to length P = 0 by the minimization objective.

The full set of constraints is listed in Model 4.5.1, and the following section explains every constraint in detail. Constraints (4.55) and (4.54) are capacity constraints and uniqueness constraints: batches have to remain within capacity b, and every job can only occupy one batch.

Constraints (4.56) define the value of P_k for every batch k as the longest p of all jobs in k. This is required in the definition of batch lateness as described above and also in (4.58), which follows the explanation above.

Constraints (4.57) ensure that no job is moved into a host-less batch, i.e. in order to move job j into batch k ($x_{jk} = 1$), job k must still be in batch k ($x_{kk} = 1$).

Constraints (4.59) implement the requirement that jobs are only moved into earlier batches.

Constraints (4.60) apply to jobs with index greater than f_L , which is the index of the job with the largest value for $L_{k,\text{single}}$. All jobs after batch f_L can be ignored in the question, based on the proof given in Appendix ??. This constraint fixes those jobs to their single-EDD batches, effectively removing them from the model. This set of constraints is rarely active but can greatly decrease the size of the model in some instances.

Constraints (4.61) through (4.63) are dominance rules implemented as lazy constraints as defined in subsection 2.2.5, and are presented in greater length in the following subsections.

4.5.1 Symmetry-breaking rule

Constraints (4.61) implement the following concept: if two jobs j_1 and j_2 are assigned as guests to batches k_1 and k_2 , and no other jobs (save for the hosts) are assigned to k_1 and k_2 ,

Min.
$$L_{\max}$$
 (4.53)
s.t. $\sum_{k} x_{jk} = 1$ $\forall j \in J$ (4.54)
 $\sum_{j} s_{j} x_{jk} \leqslant b$ $\forall k \in K$ (4.55)
 $P_{k} \geqslant p_{j} x_{jk}$ $\forall \{j \in J, k \in K | j \geqslant k\}$ (4.56)
 $x_{jk} \leqslant x_{kk}$ $\forall \{j \in J, k \in K | j \geqslant k\}$ (4.57)
 $L_{\max} \geqslant L_{k, \text{single}} - \sum_{b=0}^{k} P_{b} - p_{b}(2 - x_{bb})$ $\forall k \in K^{*}$ (4.58)
 $x_{jk} = 0$ $\forall \{j \in J, k \in K | j \geqslant k\}$ (4.59)
 $x_{jj} = 1$ $\forall \{j \in J, k \in K | j \geqslant k\}$ (4.60)
 $\forall \{j \in J, k \in K | j \geqslant k\}$ (4.61)
 $\forall \{j, j, j \in J, k \in K | j \geqslant k\}$ (4.61)
 $\forall \{j, j, j \in J, k \in K | j \geqslant k\}$ (4.61)
 $\forall \{j, j, j \in J, k \in K | j \geqslant k\}$ (4.61)
 $\forall \{j, j, j \in J, k \in K | j \geqslant k\}$ (4.62)
 $\forall \{j \in J, k \in K, j \in J, j \in J, k \in K | j \geqslant k\}$ (4.63)
 $\forall \{j \in J, k \in K, j \in J, j \in J, k \in K, j \in J$

Model 4.5.1: Move-based MIP model. Constraints marked (*) are lazy constraints (see subsection 2.2.5 for details).

then j_1 is always assigned to k_1 and j_2 to k_2 . These constraints only hold for pairs of jobs where both p_{j_1} and p_{j_2} are less than both p_{k_1} and p_{k_2} (both batches are longer than the jobs) and all possible assignments are feasible capacity-wise.

Large problem instances generate thousands of such constraints, and only very few of them will ever hold in the model. Nevertheless, they can noticeably improve solving time in some cases; declaring them as "lazy" helps keep the model small.

4.5.2 Dominance rule on "safe" moves

This applies to pairs of jobs $j_1 < j_2$ that are, potentially, competing to be guests in batch k. If both j_1 and j_2 are shorter (in terms of p) than k, we can safely enforce that in some cases, j_1 gets priority over j_2 , namely if both $p_{j_1} \ge p_{j_2}$ and $s_{j_1} \le s_{j_2}$ hold. In this case, we can state that j_2 cannot be moved into k unless j_1 has been moved somewhere also (into k or some other host batch).

Two facts allow for this set of rules:

Proposition 3. If a job j_h can either be safely moved into an earlier batch, or act as a host for later jobs $J_g = \{j_{g_1}, j_{g_2}, \ldots\}$, moving j_h is always preferable.

Proof. Let $k_{L_{\max}}$ be the batch hosting the job with maximum lateness in the final solution. If j_h is due after $k_{L_{\max}}$, moving it will have no impact on L_{\max} (see Appendix ??). If j_h is due before $k_{L_{\max}}$, but the earliest-due of J_g is due after $k_{L_{\max}}$, then moving J_g into j_h will have no impact or a worsening impact on L_{\max} . If both j_h and J_g are due before $k_{L_{\max}}$, then moving J_g will, at best, improve L_{\max} by p_{j_h} (if $\max_p(J_g) \geqslant p_{j_h}$). In this case, however, moving j_h back safely will, under any circumstances, reduce L_{\max} by p_{j_h} . The first job in J_g can host all later jobs in J_g (and possibly more, since vertical capacity was freed). The total reduction in L_{\max} in this case is improved.

Proposition 4. If a job j_1 with $d_{j_1} \leq d_{j_2}$ is both longer and thinner than j_2 ($p_{j_1} \geq p_{j_2}$, $s_{j_1} \leq s_{j_2}$), assigning it as a guest to k will produce a lower L_{max} than assigning j_2 to k.

Proof. Because of due date and size requirements, j_1 is a potential host for j_2 , Proposition 3 applies.

Generally, a longer guest job will result in a greater reduction in L_{max} , and is thus preferable. Since j_1 is thinner than j_2 , the dominance rule will not preclude (via capacity constraints) other guest jobs from being assigned to k as a consequence of prioritizing j_1 over j_2 .

4.5.3 Dominance rule on required safe moves

This constraint can be expressed logically as: if j can be safely moved into k without violating the capacity constraint, then j must be moved somewhere, or k must be empty (or both). In other words, in the context of Proposition 3, a schedule is unacceptable if job j is left in its single-EDD batch although there is room for it in an earlier batch.

The left side of the above *if-then* statement is written as $(b+1-s_j-\sum_{\substack{i=k\\i\neq j}}^{n_j}s_ix_{ik})/b$, which evaluates to 1.0 or greater iff s_k plus the sizes of guest jobs in k sum to less than $b-s_j$.

4.6 Two additional approaches tested on all models

We present two methods which, independently of the model used, lead to no change or a worsening of performance in all initial tests despite appearing very promising at first. They are mentioned here briefly only for the sake of completeness.

4.6.1 Upper bound on L_{max}

An upper bound on L_{max} can be found by using a dispatch rule to find a feasible, if not optimal, schedule. A good approach could be the "best-fit" heuristic proposed in Malapert's paper.

4.6.2 Bounding the number of batches n_{ν}

Initially, the number of batches needed is assumed to be equal to the number of jobs: $n_k = n_j$. Reducing n_k by pre-computing the maximum number of batches needed shrinks the x_{jk} matrix, and prunes potential search branches in branch-and-bound decomposition approaches.

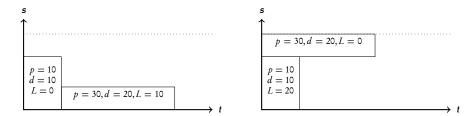


Figure 4.3: Overzealous batch elimination can increase L_{max}

Unfortunately, we cannot make a general statement that optimal solutions never have more batches than other feasible solutions – a simple counterexample is shown in figure 4.3.²

Starting out with a one-job-per-batch schedule sorted by EDD, we can explore all feasible batch configurations recursively. To generate any other feasible schedule (including the optimal solution), jobs j are rescheduled ("moved back") from their original batch k_{origin} into a prior batch k_{earlier} . This eliminates k_{origin} and requires, of course, that k_{earlier} has sufficient capacity. If the job's processing time p_j exceeds that of k_{earlier} , then the lateness of batches between k_{earlier} and k_{origin} will increase; the merit of such a move cannot be judged a priori, so it is an *unsafe* batch elimination. *Safe* eliminations, on the other hand, will never worsen L_{max} , and only they can be considered when bounding n_k a priori.

Algorithm 2 outlines a recursive method to find an upper bound on n_k , recognizing safe batch eliminations only.

This algorithm evidently requires exponential time. A relaxed variant is a possible option: if only a single unsafe move *into* a batch is possible, no safe eliminations into that

²To be more precise, we cannot state that at least one optimal solution is in the subset of feasible solutions that uses the fewest number of batches – a dominance situation that could be exploited, were it true.

Algorithm 2 Recursive algorithm to find an upper bound on n_k

```
if no open jobs left then  
update UB(n_k)
end if
find the combination of unsafe later jobs that fills up the capacity most, leaving us with capacity b_r
find all combinations x of safe later jobs that fit into b_r

repeat

x = \text{next safe job combination}
ignoreJobs \leftarrow x  
\triangleright make the moves, let the next recursion level deal with the rest of the jobs spawn and run child node with J \setminus ignoreJobs
ignoreJobs \leftarrow ignoreJobs \setminus x

until all combinations have been explored return
```

batch are considered at all and we skip to the next batch. This would greatly speed up the recursion but also significantly weaken the usefulness of the resulting upper bound.

In a branch-and-bound decomposition approach in which batches are modelled individually, an upper bound on n_k could be used to limit the depth of the search tree, or, in combination with a method to determine a lower bound on the remaining jobs' n_k at every node, to actively prune the search tree during the search. The latter method, however, would also run in exponential time as it, again, would require knapsack-type reasoning unless we use a much less powerful relaxation.

Empirical comparison of models

5.1 Experimental setup

The models were tested on a set of job lists. Malapert's paper uses benchmark job lists by Daste et al. [2008a,b], with a capacity of b = 10 and values for p_j , s_j and d_j distributed as follows:

$$p_j = U[1,99] (5.1)$$

$$s_j = U[1, 10]$$
 (5.2)

$$d_{i} = U[0,0.1] \cdot \tilde{C}_{\text{max}} + U[1,3] \cdot p_{i}$$
(5.3)

where U[a, b] is a uniform distribution between a and b, and $\tilde{C}_{\max} = \left(\sum_{j=1}^{n_j} s_j \cdot \sum_{j=1}^{n_j} p_j\right) / (b \cdot n)$ is an approximation of the time required to process all jobs.

Ten different sample job sets are used per unique value of $n_j \in \{10, 11, ..., 18, 19, 20, 25, 30, 35, 40, 45, 50\}$. The times shown in figure 5.1 are averaged over those ten instances for each n_j .

The models were run using Cplex 12.2 on an i7 Q740 CPU (1.73 GHz) in single-thread mode, with 8 GB RAM. Solving was aborted after a time of 3600 seconds (1 hour).

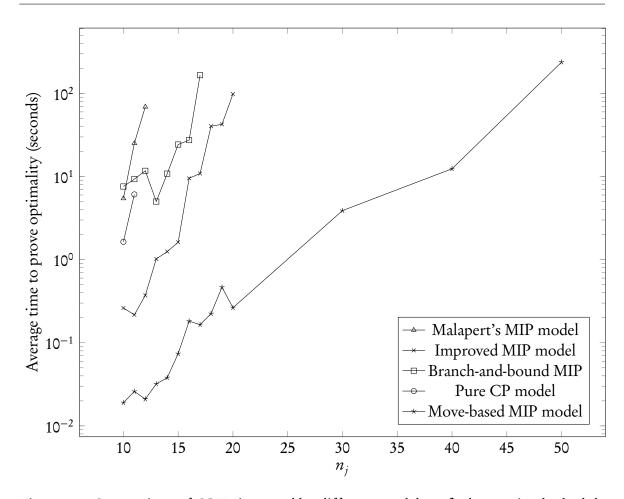


Figure 5.1: Comparison of CPU time used by different models to find an optimal schedule and prove its optimality.

5.2 Results

The CP branch-and-bound model times out on most instances and is not shown here. The pure CP model times out on one 12-job instance and gets progressively worse with more jobs; similar to Malapert's original MIP model.

The new, move-based MIP model trumps all other models' performance by at least one order of magnitude.

Unfortunately we have no such performance profile for Malapert's sequenceEDD global constraint. We do, however, have a reference in Figure 5.2, which shows the percentage of

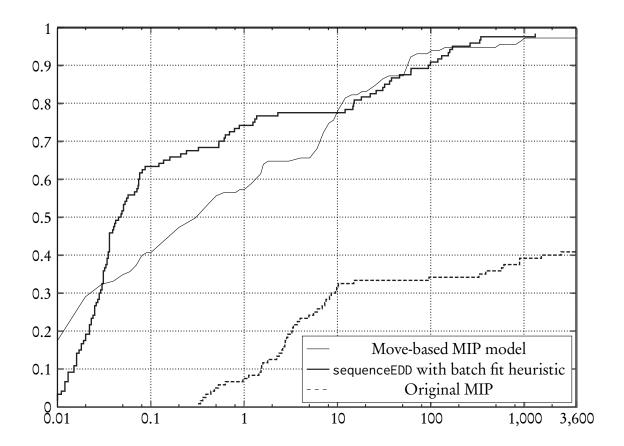


Figure 5.2: Percentage of problem instances solved in a given time

solved problem instances in a given time. The figure is based on $n_j \in \{10, 20, 50\}$, with 40 instances each (120 instances total).

5.3 Discussion

The move-based MIP model performs uses the fewest variables (columns) and constraints (rows) and performs the best. The use of lazy constraints for larger n_j is advantageous in the large majority of problem instances and thus warranted.

The improved original MIP is a great improvement on the original MIP, mostly owed to constraint (4.11) which excludes a large number of potential batch assignments.

The batch-by-batch decomposition using MIP to assemble batches, while an improvement over Malapert's MIP model, is not satisfactory; the relaxations are too lax for the

decomposition to compensate for the overhead of a naive branch-and-bound search, since we can only exclude one failed solution at a time.

5.3.1 Comparison with Malapert's sequenceEDD global constraint

Using all four filtering rules in conjunction with a batch fit search heuristic allowed Malapert to solve 38 out of 40 problem instances with $n_j = 50$ in less than an hour on a cluster of Linux machines, each node of which using 48 GB of RAM and two quad core 2.4 GHz processors (that is, using considerably more powerful hardware than was used for this thesis).

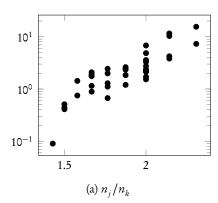
The move-based MIP model is able to achieve a comparable performance (39 of 40 instances solved within an hour). While it is, on average, slower on problems with low n_j (compare Figure 5.2), it is able to prove optimality quicker on most instances that take longer than 10 seconds to solve—ostensibly the more critical metric.

5.4 Further performance tests

A better understanding of the relationships between model performance and problem instance characteristics can help in the development of faster models. This is especially true with regard to redundant and/or lazy constraints that are used only to achieve a tighter formulation at the price of a larger model: using such constraints is often warranted only for particularly challenging problem instances. The following paragraphs outline some such considerations.

5.4.1 Correlation between disjunctivity and solving time

The mean number of jobs per batch (n_j/n_k) in the optimal solution is strongly correlated with the time needed to solve to optimality. This is unsurprising given that a problem with a great number of large jobs (i.e. large s_j) will be of relatively low difficulty in terms of bin packing. This is illustrated by Figure 5.3a. An excellent predictor of this correlation appears to be what Baptiste and Le Pape Baptiste and Le Pape [2000] call disjunction ratio:



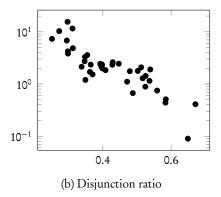


Figure 5.3: Correlation between time to prove optimality in seconds (ordinates) and measures of disjunctivity (abscissae), based on 40 instances with $n_j = 30$. Subfigure (a) shows the average number of jobs per batch in the optimal solution. Subfigure (b) shows the *disjunction ratio* as defined in Section 5.4.1.

the ratio between the number of job pairs which cannot run in parallel (as $s_{j_1} + s_{j_2} > b$) and the total number of job pairs n_j^2 . The relationship is shown in Figure ??.

As a corrolary to the above, the solving time is greatly dependent on the capacity of the machine. Larger capacities, relative to the average job size \bar{s}_j , will cause longer solving times as the disjunction ratio decreases.

5.4.2 Uniform distribution of due dates

The due dates in the test instances used by Malapert and in the previous tests were generated by a fairly complex formula designed such that jobs are likely to fall into *buckets*, i.e. larger sets of jobs with equal due dates.

Malapert's sequenceEDD global constraint exploits the notion of buckets extensively, and would likely perform worse on instances that exhibit no bucketing at all (though we could not test this for lack of a working implementation).

Our move-based model uses only the latest-due batch in every bucket to define the value of $L_{\rm max}$. As a result, the size of the model is directly dependent on the number of buckets. We were, however, unable to find evidence of such a relationship: Figure 5.4 shows the

¹The term originates from the dichotomy between *disjunctive resources* which can only process one job at a time, and *(highly) cumulative resources* which can process many jobs in parallel.

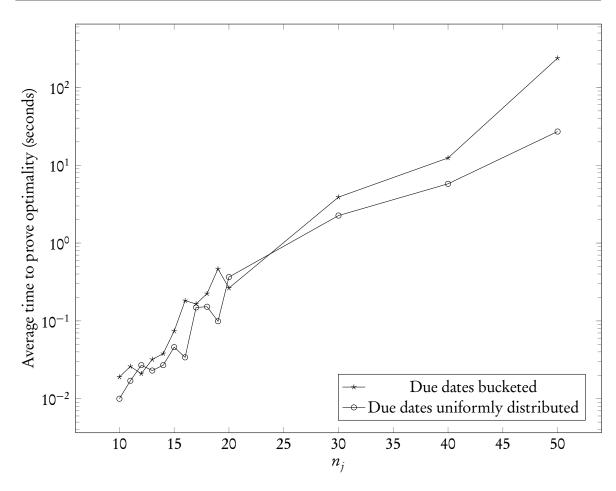


Figure 5.4: Comparison of time to prove optimality required by the move-based MIP model depending on the distribution of due dates.

performance of the move-based MIP on non-bucketed ($d_j = U[1,99]$) instances compared to bucketed instances.

Indeed, average performance improves with uniformly distributed due dates on instances with larger n_j .

Conclusion

The problem of scheduling jobs of non-identical sizes and processing times on a batch resource with a limited capacity arises frequently in the context of furnaces, autoclaves, driers, or other machinery where one particular job determines the processing time of the entire batch. The problem of minimizing the maximum lateness is given in situations in which all customers (or downstream processes) are to be treated as fairly as possible.

The problem was approached by Daste Daste et al. [2008a] using a branch-and-price model and by Malapert Malapert et al. [2012] using a new sequenceEDD global constraint. In this paper we have introduced a set of improvements to Malapert's basic MIP model, a simple CP model, a decomposition method, and a new MIP based on the idea of independent search moves expressed in a single formulation.

The new move-based MIP model rivals the performance of Malapert's sequenceEDD constraint particularly with larger problem instances ($n_j > 30$). The great simplicity of our formulation allows production managers to integrate it into existing models without having to implement highly problem-specific filtering rules. Furthermore, it is easily extended with additional redundant or lazy constraints if necessary.

6.1 Future Work

All models described in this paper have weaknesses. The improved original MIP model lacks the fine-tuning and lazy constraints enjoyed by the move-based MIP. The CP model could likely benefit from additional redundant constraints. The decomposition approach

needs a better pruning method. The move-based MIP's lazy constraints could certainly be fine-tuned based on a more generous definition of "safe" moves and conditional activation relying on the problem's *disjunction ratio*. Such potential improvements are described in Appendix ??.

6.1.1 Application to related problems

The promising computational performance of the move-based MIP model might find application in similar problems, e.g. completion-time minimization on batch machines (which is equivalent to L_{\max} minimization with equal due dates), or L_{\max} minimization with release dates (i.e. coupling x_{jk} to batch start times, which can be expressed using an exact analogue to constraint (4.58)).

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Appendix: Potential improvements to the models

The following section details some redundant constraints, tighter variable bounds and potential fine-tuning opportunities that were not implemented in this papier, but may be of use in further research.

A.1 CP Model

A.1.1 Constraint on number of batches with length $P_k > p$

Since batches take on the processing time of their longest job, there is at least one batch with $P = \max_{j} p_{j}$. We can proceed to fill batches with jobs, ordered by non-increasing processing time, based on algorithm 3.

At the end of this algorithm, we can state:

$$\mbox{globalCardinality}(\{1,\dots,n_{j}\},\{P_{k,\min},\dots,P_{k-1,\min}-1\},P_{k}) \quad \forall k \in \{k_{0},\dots,k_{LB(n_{k})}\}, \tag{A.1}$$

where the constraint takes three arguments (cards, vals and vars) and $P_{k,\min}$ denotes the minimum length of batch k.

The algorithm sorts jobs by non-increasing *p*, and then fills batches job by job. If a job fits into a previous batch, it is assigned there. If a job fits into multiple previous batches, it is assigned to the batch with the smallest remaining capacity. This is called *best-fit decreasing*

Algorithm 3 Generating lower bounds on batch lengths

```
▶ initialize all jobs as unassigned jobs
n_k \leftarrow 1; S_k \leftarrow \{0\}; P_{k, \min} \leftarrow \{0\}
                                                                                        ▶ Create one empty batch of size and length zero
sort J^* by processing time, non-increasing
                                                                                               ▶ select job for assignment, longest job first
     i \leftarrow I^*.pop()
    loop through all n_k existing batches k, first batch first
                                                                                                                               ▶ no feasible batch
         c_{\min} = b
                                                                                        > currently known minimum remaining capacity
         if s_j < b - S_k and b - S_k < c_{\min} then
             k_p \leftarrow k_p; c_{\min} \leftarrow b - S_k
    end loop
    if |k_n| = 1 then
                                                                                                                      \triangleright assign job j to batch k_p
         S_{k_p} \leftarrow S_{k_p} + s_j
         if n_k < LB(n_k) then
              n_k \leftarrow n_k + 1
                                                                                                                                ▶ open new batch
             S_{n_k} \leftarrow s_j; P_{n_k, \min} \leftarrow p_j
                                                                                                          \triangleright assign s_i and p_i to the new batch
              leave the loop now and end.
         end if
    end if
until J^* is empty
```

rule, and works as follows: let J^* be the set of jobs sorted by p, then at least one batch will be as long as the longest job j_1^* . If the next n jobs fit into this batch, then there is at least one batch not shorter than j_{n+1}^* , and similarly for subsequent batches.

Unfortunately, the optimal solution may perform better than the packing heuristic in terms of "vertical" (s_j) bin packing, and may thus require fewer batches. We therefore need to find a lower bound $LB(n_k)$ on the number of batches, and we can only guarantee the first $LB(n_k)$ of the above constraints to hold in the optimal solution. Finding a true lower bound is a two-dimensional bin packing problem, which is NP-hard. A possible but naive lower bound is j_0 , the number of jobs, ordered by decreasing s_j , that can never fit into a batch together.

A.1.2 Constraint on number of batches with length $D_k > d$

In a similar fashion, we can determine that the second batch must be due no later than the earliest-due job j_{m+1} that can *not* fit into the first batch – if we sort jobs by due date and fit

the earliest m jobs into the first batch – and so on for subsequent batches. Once again, since best-fit decreasing may not perform optimally in terms of "vertical" packing, this may not be valid for batches beyond the known $LB(n_k)$.

A.1.3 All-different constraints on P_k and D_k

Furthermore, if all jobs have different processing times, all batches will have different processing times as well: $\operatorname{alldifferent}(P_k)$. If m out of n_j jobs have different processing times, we can still enforce k_alldifferent(P_k , m). Some work on k_alldifferent constraints has been done in [Lardeux et al.].

Similarly, we knows that the constraint k_alldifferent(D_k , m) must be true if m out of n_i jobs have different due dates.

A.2 Decomposition approaches

A.2.1 Potential heuristics

Improve the initial $L_{\text{max,incumbent}}$ A better initial upper bound on L_{max} can help prune some branches of the search tree from the outset. There are several dispatch rules (or maybe other heuristics?) that could be explored to do this better.

Improve $L_{\text{max,incumbent}}$ during search It may be useful to use a heuristic like above to "complete the schedule" once a promising partial schedule has been generated. I have yet to identify situations where this is always helpful.

A.3 Move-based MIP model

A.3.1 Conditional activation of lazy constraints

The disjunction ratio measure introduced by? may be useful in conditionally activating or omitting certain lazy constraint sets, as it is known that they often have a net slowing effect on the solver when dealing with less difficult problems.

A.3.2 Improvement of "safe" move definition

The concept of "safe" moves is used in several constraints. It describes moves in which a job j is moved into an earlier batch k and $p_j \leq p_k$.

The definition of "safe" can be extended to guest jobs j that fulfill the following condition: $p_j \leq LB(L_{\max}) + d_k - UB(S_k)$, where $LB(L_{\max})$ is a lower bound on the optimal value of L_{\max} and $UB(S_k)$ is an upper bound on the start time of batch k. The term represents the longest processing time that batch k can be extended to without violating a known lower bound on the maximum lateness.

The a-priori computation of upper bounds on batch start times is challenging, and can only be based on capacity violations ("what is the longest combination of guest jobs that can be placed into batches up to and including k?"), which will likely make for relatively weak bounds. Nevertheless, a stronger notion of "safe" may be beneficial in some instances.

A.4 Move-back search

The following is a proof referenced in Proposition 3, but it could easily be turned into a standalone local search method that finds good solutions quickly.

Proposition 5. Begin with any EDD schedule (such as the single-EDD schedule). Moving any job j into an earlier batch k will never shift the position of L_{max} to the right (into a later bucket).

Proof. In any partial schedule following EDD, let k be the batch with maximum lateness $L_k = L_{\max}$. If multiple batches have equal lateness L_{\max} , let L_k refer to the latest-due of these batches in the schedule. It has processing time p_k . Then the lateness of the batch before k must be $L_{k-1} \geqslant L_k - p_k$ as a consequence of the EDD sequencing. Any batch following k can have a lateness no greater than $L_k - 1$.

- Moving any single job from a batch following k into a batch before k will worsen L_{\max} , but not change its position. Such a move is never necessary to arrive at an optimal solution.

- Moving back any single job from a batch before k safely¹ will improve L_{max} , but not change its position.
- Moving back a single pre-k job j from a batch β into an earlier batch α unsafely will reduce L_k by p_{α} ; L_{\max} may still be in k, or it may be found in any batch between (and including) α and β , since their lateness $L_{[\alpha,\dots,\beta]}$ is now increased by $p_j p_{\alpha}$.
- If the max-lateness job j is single itself, it can be moved from k into an earlier batch α . If this is done *safely*, the batch immediately preceding k will still have a lateness $L_{k-1} \ge L_k p_j$. All batches after k will have their lateness reduced by p_j , but since their maximum lateness did not exceed $L_k 1$ before the move, it will now be at least 1 less than that of batch k-1.
- If the max-lateness job j was moved back *unsafely* into a batch α , batches between and including α and k-1 now have their lateness increased by p_j-p_α , while batches after k have their lateness decreased by p_α . Again, batch L_{k-1} would exceed the maximum lateness of batches after k.

This shows that after a sequence of operations in which single jobs are moved into earlier batches, $L_{\rm max}$ will never shift to the right. In fact, this means that all jobs after the $L_{\rm max}$ -job in a single-job EDD schedule can be ignored in the scheduling problem entirely, although this turns out to be quite inconsequential as that job is often at or near the end of the single-job EDD schedule, resulting from the fact that $L_{k-1} \geqslant L_k - p_k$. By the same token, high-quality solutions often have their $L_{\rm max}$ in an early batch. This may give rise to some sort of decomposition method in which jobs are strategically moved back, and where we are trying to move $L_{\rm max}$ to the left as m

¹ for the meaning of "safe" and "unsafe" moves, compare section 4.6.2.

Appendix: Other approaches

Appendix: Result tables

C.1 Hello. This is the first part of the first Appendix Blabla.

Appendix: Code listings